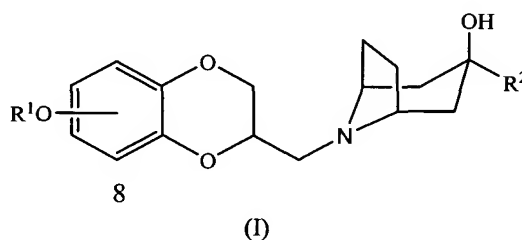


This listing of claims will replace all prior versions, and listings, of claims in the application.

Listing of Claims:

Claims 1 to 25 (*cancelled*)

26. (*currently amended*) A method of treating a subject suffering from a condition selected from the group consisting of neurodegenerative disease, eating disorders, disorders of thermoregulation, sleep dysfunction and sexual dysfunction, comprising the step of: which comprises providing to the subject suffering from said condition, a therapeutically effective amount of a compound of formula I



wherein

R¹ is a straight-chained alkyl of 1 to 6 carbon atoms, or a branched chain alkyl of 3 to 8 carbon atoms; and

R² is phenyl, naphthyl, anthracyl, phenanthryl, pyridyl, pyrimidyl, triazinyl, furyl, pyrrolyl, pyrazolyl, indolyl, imidazolyl, benzofuryl, benzothienyl, oxazolyl, or thiazolyl each optionally substituted with 0 to 3 substituents selected from straight-chain alkyl of 1 to 6 carbon atoms, branched-chain alkyl of 3 to 8 carbon atoms, alkoxy of 1 to 6 carbon atoms, mono- or dialkylamino of 1 to 6 carbon atoms, nitro, halo, amino, cyano, trifluoromethyl, trifluoromethoxy and hydroxy;

and or a pharmaceutically acceptable salts salt thereof.

Claims 27 to 32 (*cancelled*)

33. (*new*) A method according to claim 26, wherein said subject is a human.

34. *(new)* A method according to claim 26, wherein R^1 is a straight-chained alkyl of 1 to 3 carbon atoms, or a branched chain alkyl of 3 to 6 carbon atoms.
35. *(new)* A method according to claim 26, wherein R^1 is a straight-chained alkyl of 1 or 2 carbon atoms.
36. *(new)* A method according to claim 26, wherein R^2 is phenyl, naphthyl, pyridyl, pyrimidyl, furyl, pyrrolyl, pyrazolyl, indolyl, imidazolyl, benzofuryl, or benzothienyl; each optionally substituted with 1 to 3 substituents the same or different selected from straight-chain alkyl of 1 to 3 carbon atoms, branched-chain alkyl of 3 to 6 carbon atoms, alkoxy of 1 to 3 carbon atoms, mono- or di-alkylamino in which each alkyl group has 1 to 3 carbon atoms, nitro, amino, cyano, halogen, trifluoromethyl, trifluoromethoxy, and hydroxy.
37. *(new)* A method according to claim 26, wherein R^2 is phenyl, naphthyl, pyridyl, pyrrolyl, indolyl, or benzothienyl; each optionally substituted with 1 to 3 substituents the same or different selected from nitro, amino, cyano, halogen, trifluoromethyl, trifluoromethoxy, and hydroxy.
38. *(new)* A method according to claim 26, wherein R^2 is trifluoromethylphenyl or methoxyphenyl.
39. *(new)* A method according to claim 26, wherein the R^1O substituent is bonded to the 1,4-benzodioxan nucleus at the 8 position.
40. *(new)* A method according to claim 26, wherein R^1 is a straight-chained alkyl of 1 to 3 carbon atoms, or a branched chain alkyl of 3 to 6 carbon atoms and R^2 is phenyl, naphthyl, pyridyl, pyrimidyl, furyl, pyrrolyl, pyrazolyl, indolyl, imidazolyl, benzofuryl, or benzothienyl; each optionally substituted with 0 to 3 substituents

selected from straight-chain alkyl of 1 to 3 carbon atoms, branched-chain alkyl of 3 to 6 carbon atoms, alkoxy of 1 to 3 carbon atoms, mono- or di-alkylamino in which each alkyl group has 1 to 3 carbon atoms, halogen, trifluoromethyl, trifluoromethoxy, and hydroxy.

41. *(new)* A method according to claim 26, wherein R¹ is a straight-chained alkyl of 1 or 2 carbon atoms, and R² is phenyl, naphthyl, pyridyl, pyrrolyl, indolyl, or benzothienyl; each optionally substituted with a 0 to 3 substituents selected from nitro, amino, cyano, halogen, trifluoromethyl, trifluoromethoxy, and hydroxy.
42. *(new)* A method according to claim 26, wherein R¹ is a straight chain alkyl of 1 or 2 carbon atoms and R² is trifluoromethylphenyl or methoxyphenyl.
43. *(new)* A method according to claim 26, wherein said compound is (S)-8-(8-ethoxy-2,3-dihydrobenzo-[1,4]dioxin-2-ylmethyl)-3-naphthalen-2-yl-8-aza-bicyclo[3.2.1]octan-3-ol or a pharmaceutically acceptable salt thereof.
43. *(new)* A method according to claim 26, wherein said compound is (S)-8-(8-ethoxy-2,3-dihydro-benzo[1,4]dioxin-2-ylmethyl)-3-phenyl-8-aza-bicyclo[3.2.1]octan-3-ol or a pharmaceutically acceptable salt thereof.
44. *(new)* A method according to claim 26, wherein said compound is (S)-3-benzo[b]thiophen-3-yl-8-(8-ethoxy-2,3-dihydro-benzo[1,4]dioxin-2-ylmethyl)-8-aza-bicyclo[3.2.1]octan-3-ol or a pharmaceutically acceptable salt thereof.
45. *(new)* A method according to claim 26, wherein said compound is 8-{[(2S)-8-ethoxy-2,3-dihydrobenzo-[1,4]dioxin-2-yl]methyl)-3-pyridin-2-yl-8-aza-bicyclo[3.2.1]octan-3-ol or a pharmaceutically acceptable salt thereof.

46. *(new)* A method according to claim 26, wherein said compound is 8-{[(2S)-8-ethoxy-2,3-dihydrobenzo-[1,4]dioxin-2-yl)methyl]-3-(3-trifluoromethyl-phenyl)-8-aza-bicyclo[3.2.1]octan-3-ol or a pharmaceutically acceptable salt thereof.
47. *(new)* A method according to claim 26, wherein said compound is 8-{[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl)methyl]-3-(2-methoxyphenyl)-8-azabicyclo[3.2.1]octan-3-ol or a pharmaceutically acceptable salt thereof.
48. *(new)* A method according to claim 26, wherein said compound is 8-{[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl)methyl]-3-[3-(trifluoromethyl)phenyl]-8-azabicyclo[3.2.1]octan-3-ol or a pharmaceutically acceptable salt thereof.
49. *(new)* A method according to claim 26, wherein said compound is 8-{[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl)methyl]-3-(2-pyridinyl)-8-azabicyclo[3.2.1]octan-3-ol or a pharmaceutically acceptable salt thereof.
50. *(new)* A method according to claim 26, wherein said compound is 3-(1-benzothien-3-yl)-8-{[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl)methyl]-8-azabicyclo[3.2.1]octan-3-ol or a pharmaceutically acceptable salt thereof.
51. *(new)* A method according to claim 26, wherein said compound is 8-{[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl)methyl]-3-phenyl-8-azabicyclo[3.2.1]octan-3-ol or a pharmaceutically acceptable salt thereof.
52. *(new)* A method according to claim 26, wherein said compound is 3-((2S)-8-methoxy-2,3-dihydrobenzo-[1,4]dioxin-2-yl)methyl)-8-naphthalen-2-yl-3-azabicyclo[3.2.1]octan-8-ol or a pharmaceutically acceptable salt thereof.